

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034527.D
 Acq On : 10 Oct 2024 18:43
 Operator : RC/JU
 Sample : P4263-09
 Misc :
 ALS Vial : 103 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EB04-20241001

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/11/2024
 Supervised By :mohammad ahmed 10/14/2024

Quant Time: Oct 11 00:22:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 00:21:08 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	4333	0.400	ng	0.00
7) Naphthalene-d8	10.127	136	11948	0.400	ng	0.00
13) Acenaphthene-d10	14.019	164	6655	0.400	ng	0.00
19) Phenanthrene-d10	16.783	188	12408	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	8167	0.400	ng	0.00
35) Perylene-d12	23.119	264	10036	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1514	0.123	ng	0.00
5) Phenol-d6	6.585	99	775m	0.054	ng	0.01
8) Nitrobenzene-d5	8.525	82	2362	0.258	ng	0.01
11) 2-Methylnaphthalene-d10	11.731	152	5639	0.320	ng	0.00
14) 2,4,6-Tribromophenol	15.542	330	822	0.273	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	8013	0.296	ng	0.00
27) Fluoranthene-d10	18.829	212	12868	0.411	ng	0.00
31) Terphenyl-d14	19.452	244	8971	0.550	ng	0.00
Target Compounds						
9) Naphthalene	10.180	128	809	0.025	ng	# 69

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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